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data from INPADOC  
NEWS 4 FEB 28 BABS - Current-awareness alerts (SDIs) available  
NEWS 5 MAR 02 GBFULL: New full-text patent database on STN  
NEWS 6 MAR 03 REGISTRY/ZREGISTRY - Sequence annotations enhanced  
NEWS 7 MAR 03 MEDLINE file segment of TOXCENTER reloaded  
NEWS 8 MAR 22 KOREAPAT now updated monthly; patent information enhanced  
NEWS 9 MAR 22 Original IDE display format returns to REGISTRY/ZREGISTRY  
NEWS 10 MAR 22 PATDPASPC - New patent database available  
NEWS 11 MAR 22 REGISTRY/ZREGISTRY enhanced with experimental property tags  
NEWS 12 APR 04 EPFULL enhanced with additional patent information and new  
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NEWS 13 APR 04 EMBASE - Database reloaded and enhanced  
NEWS 14 APR 18 New CAS Information Use Policies available online  
NEWS 15 APR 25 Patent searching, including current-awareness alerts (SDIs),  
based on application date in CA/CAPLUS and USPATFULL/USPAT2  
may be affected by a change in filing date for U.S.  
applications.  
NEWS 16 APR 28 Improved searching of U.S. Patent Classifications for  
U.S. patent records in CA/CAPLUS  
NEWS 17 MAY 23 GBFULL enhanced with patent drawing images  
NEWS 18 MAY 23 REGISTRY has been enhanced with source information from  
CHEMCATS  
NEWS 19 JUN 06 The Analysis Edition of STN Express with Discover!  
(Version 8.0 for Windows) now available  
NEWS 20 JUN 13 RUSSIAPAT: New full-text patent database on STN  
NEWS 21 JUN 13 FRFULL enhanced with patent drawing images  
NEWS 22 JUN 27 MARPAT displays enhanced with expanded G-group definitions  
and text labels  
NEWS 23 JUL 01 MEDICONF removed from STN  
NEWS 24 JUL 07 STN Patent Forums to be held in July 2005  
NEWS 25 JUL 13 SCISEARCH reloaded  
NEWS 26 JUL 20 Powerful new interactive analysis and visualization software,  
STN AnaVist, now available  
NEWS 27 AUG 11 Derwent World Patents Index(R) web-based training during  
August  
NEWS 28 AUG 11 STN AnaVist workshops to be held in North America  
  
NEWS EXPRESS JUNE 13 CURRENT WINDOWS VERSION IS V8.0, CURRENT  
MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),  
AND CURRENT DISCOVER FILE IS DATED 13 JUNE 2005

08/30/2005 10626155.trn

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NEWS INTER	General Internet Information
NEWS LOGIN	Welcome Banner and News Items
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\* \* \* \* \* STN Columbus \* \* \* \* \*

FILE 'HOME' ENTERED AT 08:36:59 ON 30 AUG 2005

=>

Uploading

THIS COMMAND NOT AVAILABLE IN THE CURRENT FILE

Do you want to switch to the Registry File?

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Switching to the Registry File...

Some commands only work in certain files. For example, the EXPAND command can only be used to look at the index in a file which has an index. Enter "HELP COMMANDS" at an arrow prompt (=>) for a list of commands which can be used in this file.

=> FILE REGISTRY

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'REGISTRY' ENTERED AT 08:37:12 ON 30 AUG 2005

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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 29 AUG 2005 HIGHEST RN 862072-85-3

DICTIONARY FILE UPDATES: 29 AUG 2005 HIGHEST RN 862072-85-3

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JULY 14, 2005

Please note that search-term pricing does apply when conducting SmartSELECT searches.

\*\*\*\*\*  
\*  
\* The CA roles and document type information have been removed from \*  
\* the IDE default display format and the ED field has been added, \*

08/30/2005 10626155.trn

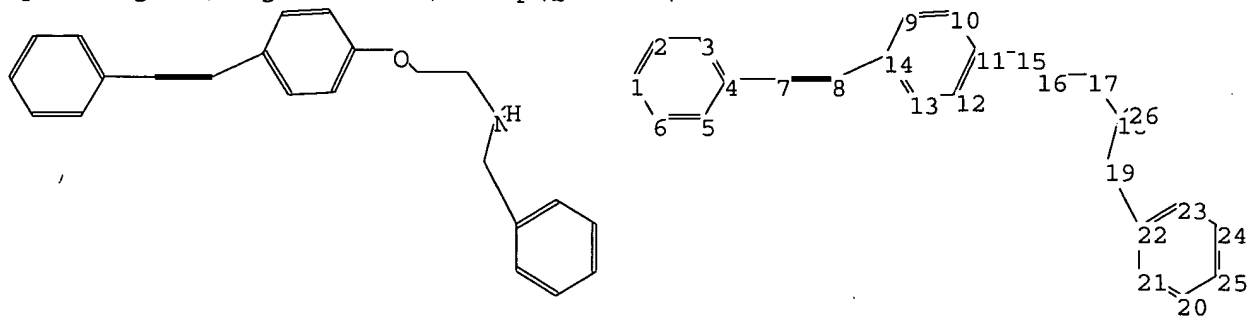
\* effective March 20, 2005. A new display format, IDERL, is now \*  
\* available and contains the CA role and document type information. \*  
\*  
\*\*\*\*\*

Structure search iteration limits have been increased. See HELP SLIMITS  
for details.

Experimental and calculated property data are now available. For more  
information enter HELP PROP at an arrow prompt in the file or refer  
to the file summary sheet on the web at:  
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>

Uploading C:\Program Files\Stnexp\Queries\10626155.str



chain nodes :

7 8 15 16 17 18 19 26

ring nodes :

1 2 3 4 5 6 9 10 11 12 13 14 20 21 22 23 24 25

chain bonds :

4-7 7-8 8-14 11-15 15-16 16-17 17-18 18-19 18-26 19-22

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 9-10 9-14 10-11 11-12 12-13 13-14 20-21 20-25

21-22 22-23 23-24 24-25

exact/norm bonds :

11-15 15-16 17-18 18-19

exact bonds :

4-7 7-8 8-14 16-17 18-26 19-22

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6 9-10 9-14 10-11 11-12 12-13 13-14 20-21 20-25

21-22 22-23 23-24 24-25

isolated ring systems :

containing 1 : 9 : 20 :

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:Atom 10:Atom

11:Atom 12:Atom 13:Atom 14:Atom 15:CLASS 16:CLASS 17:CLASS 18:CLASS

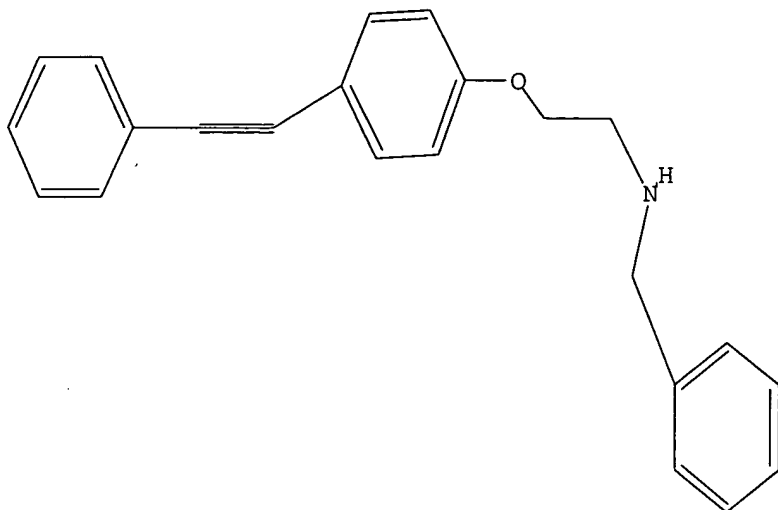
19:CLASS 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom 26:CLASS

L1 STRUCTURE UPLOADED

=> d l1

L1 HAS NO ANSWERS

L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> s l1

SAMPLE SEARCH INITIATED 08:37:29 FILE 'REGISTRY'  
SAMPLE SCREEN SEARCH COMPLETED - 9 TO ITERATE

100.0% PROCESSED 9 ITERATIONS 0 ANSWERS  
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
PROJECTED ITERATIONS: 9 TO 360  
PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=> s l1 sss full

FULL SEARCH INITIATED 08:37:36 FILE 'REGISTRY'  
FULL SCREEN SEARCH COMPLETED - 327 TO ITERATE

100.0% PROCESSED 327 ITERATIONS  
SEARCH TIME: 00.00.01

37 ANSWERS

L3 37 SEA SSS FUL L1

=> FIL HCAPLUS

COST IN U.S. DOLLARS

FULL ESTIMATED COST

SINCE FILE	TOTAL
ENTRY	SESSION
161.33	161.54

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FILE 'HCAPLUS' ENTERED AT 08:37:44 ON 30 AUG 2005  
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FILE COVERS 1907 - 30 Aug 2005 VOL 143 ISS 10  
FILE LAST UPDATED: 29 Aug 2005 (20050829/ED)

New CAS Information Use Policies, enter HELP USAGETERMS for details.

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l3  
L4

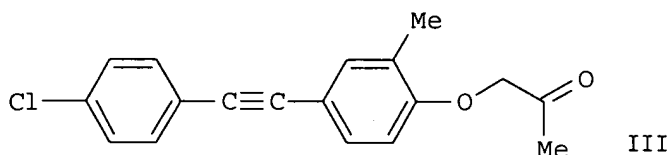
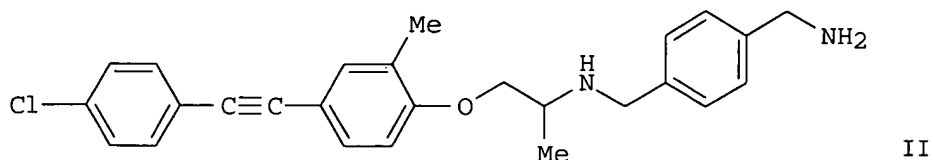
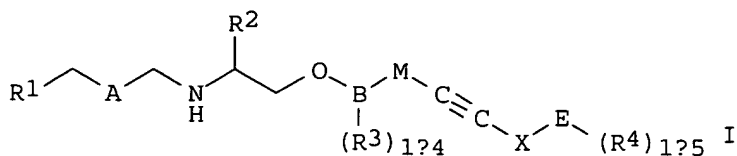
4 L3

=> d l4 ibib abs hitstr tot

L4 ANSWER 1 OF 4 HCAPLUS COPYRIGHT 2005 ACS on STN  
ACCESSION NUMBER: 2005:78301 HCAPLUS  
DOCUMENT NUMBER: 142:176433  
TITLE: A preparation of acetylenic compounds, useful in the treatment of inflammatory disorders  
INVENTOR(S): Beers, Scott; Malloy, Elizabeth A.; Wachter, Michael P.  
PATENT ASSIGNEE(S): USA  
SOURCE: U.S. Pat. Appl. Publ., 19 pp.  
CODEN: USXXCO  
DOCUMENT TYPE: Patent  
LANGUAGE: English  
FAMILY ACC. NUM. COUNT: 1  
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 2005020838	A1	20050127	US 2003-626155	20030724
PRIORITY APPLN. INFO.:			US 2003-626155	20030724
OTHER SOURCE(S):	MARPAT	142:176433		

GI



AB The invention relates to a preparation of acetylenic compds. of formula I [wherein: A is cycloalkyldiyl, cyclic heteroalkyldiyl, or (hetero)aryldiyl; B is (hetero)aryldiyl; E is (hetero)aryldiyl; M and X are independently (CH<sub>2</sub>)<sub>0-4</sub>; R<sub>1</sub> is cycloalkyl, cyclic heteroalkyl, or (hetero)aryl, etc.; R<sub>2</sub> is H, alkyl, alkoxy, CHO, CO<sub>2</sub>H, or NH<sub>2</sub>, etc.; R<sub>3</sub> and R<sub>4</sub> are independently H, alkyl, CHO, cycloalkyl, or aryl, etc.], useful in the treatment of inflammatory disorders. For instance, acetylene derivative II (oxidase inhibition: IC<sub>50</sub> = 0.7 μM) was prepared via reductive amination of ketone III by 1,4-bis(aminomethyl)benzene.

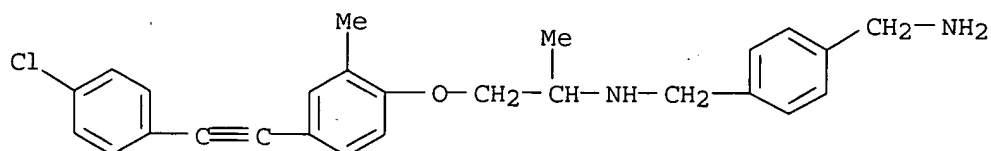
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651330-49-3P 651330-50-6P 651330-51-7P  
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651330-55-1P 651330-58-4P 651330-59-5P  
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651330-65-3P 651330-68-6P 773847-69-1P  
776293-45-9P 792905-81-8P 832116-38-8P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of acetylenic compound useful in the treatment of inflammatory disorders)

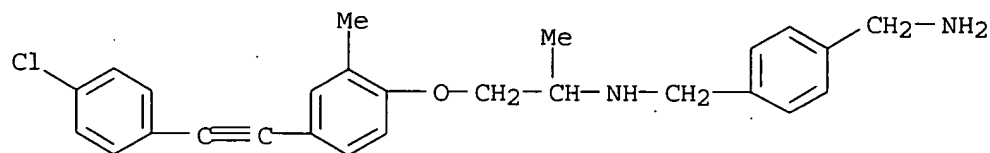
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RN 651330-37-9 HCAPLUS

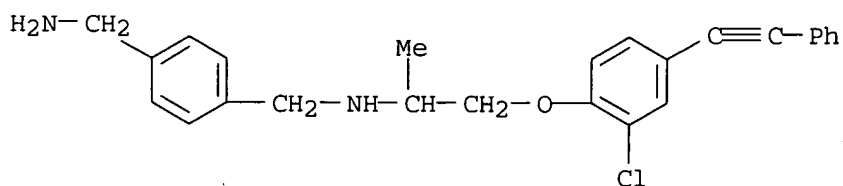
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● 2 HCl

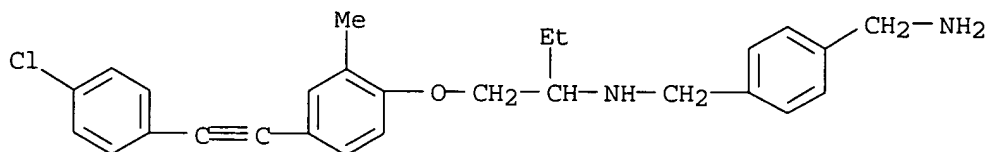
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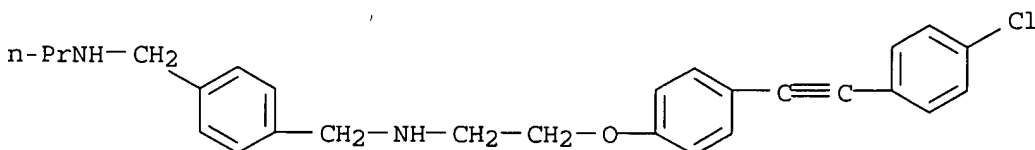
RN 651330-39-1 HCAPLUS

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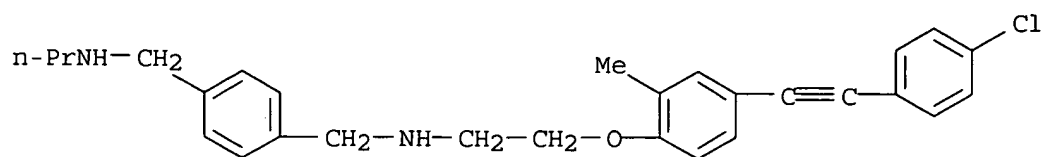
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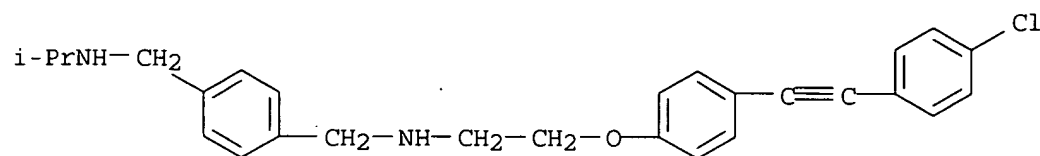
RN 651330-45-9 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-propyl- (9CI) (CA INDEX NAME)



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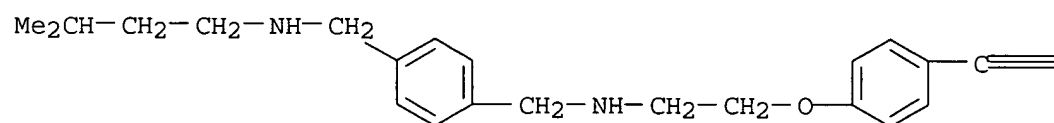
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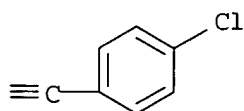
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CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]phenoxy]ethyl]-N'-(3-methylbutyl)- (9CI) (CA INDEX NAME)

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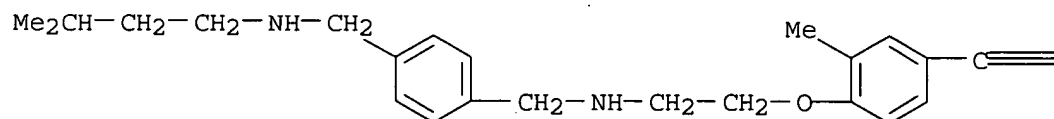
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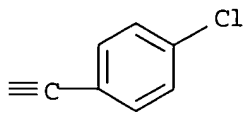
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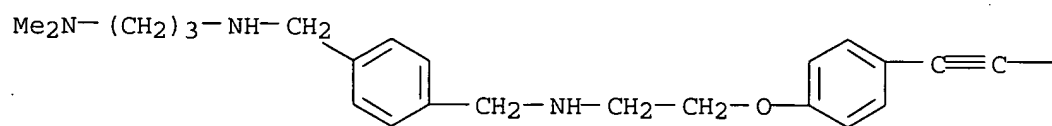
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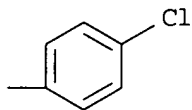
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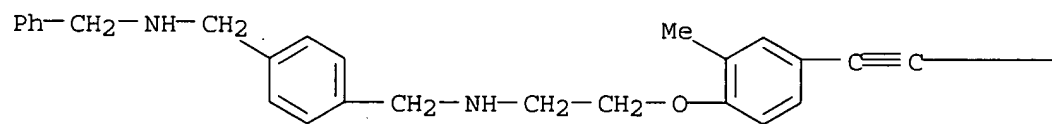
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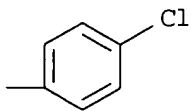
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PAGE 1-A



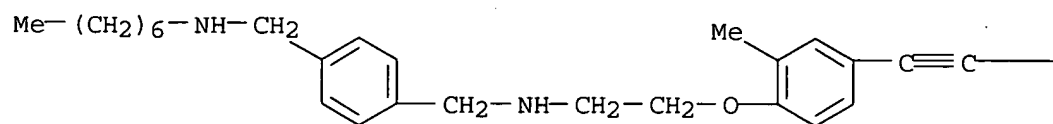
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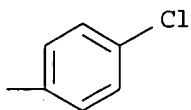
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PAGE 1-A



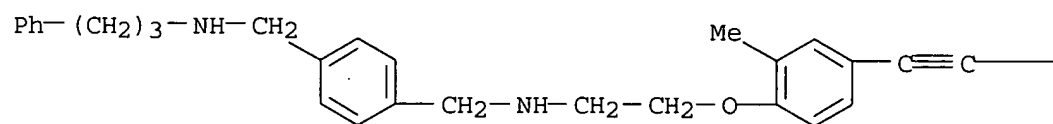
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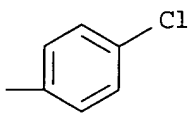
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PAGE 1-A



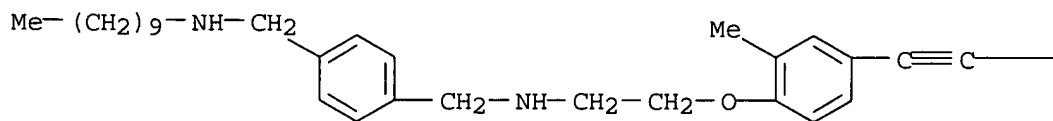
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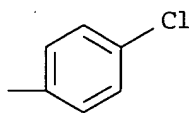
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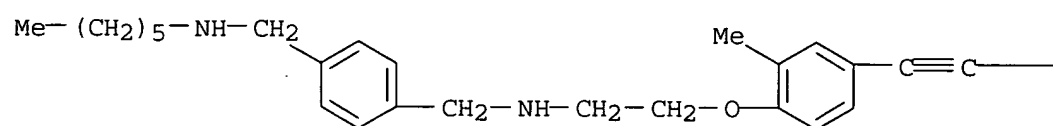
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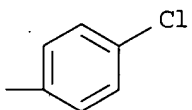
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PAGE 1-A



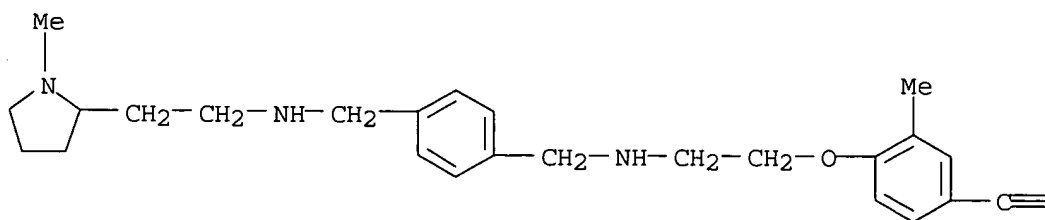
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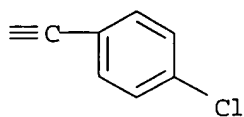
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CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-[2-(1-methyl-2-pyrrolidinyl)ethyl]- (9CI) (CA INDEX NAME)

PAGE 1-A

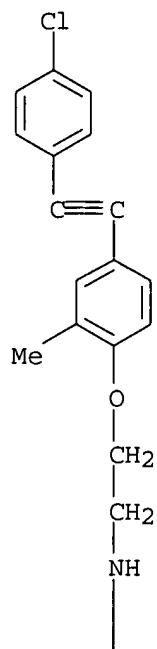


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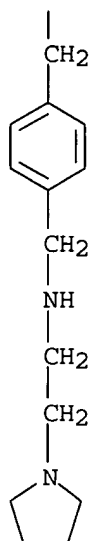


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PAGE 1-A



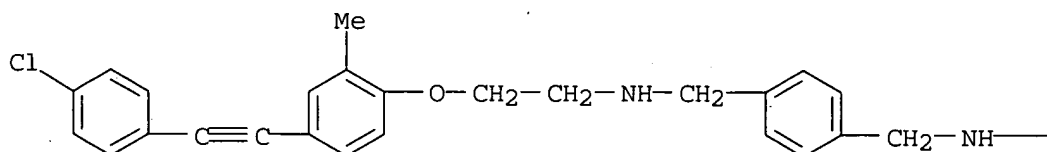
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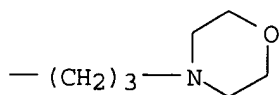
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PAGE 1-A



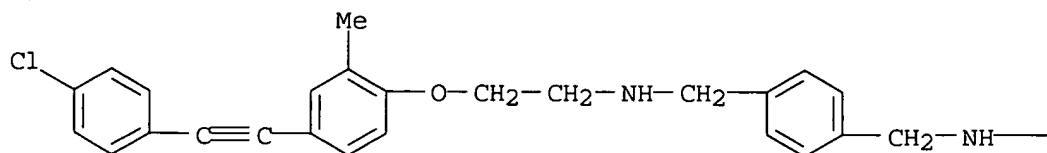
PAGE 1-B



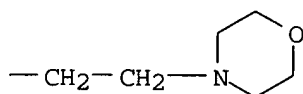
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PAGE 1-A

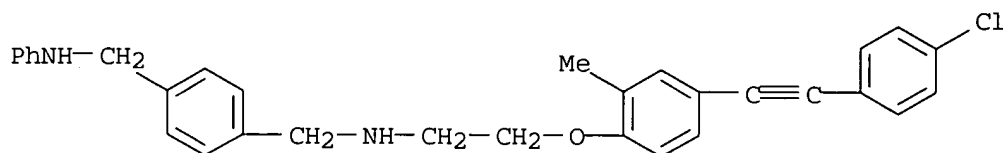


PAGE 1-B



RN 651330-63-1 HCAPLUS

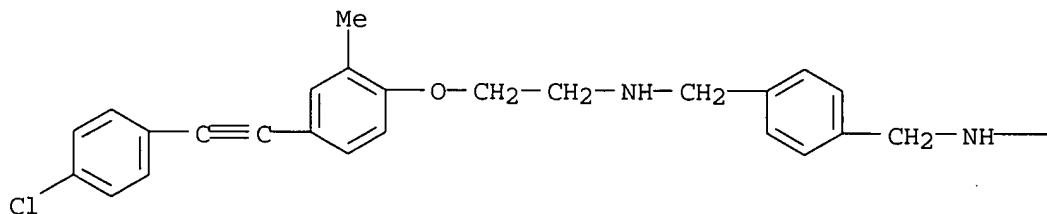
CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-phenyl- (9CI) (CA INDEX NAME)



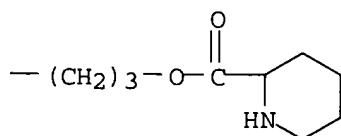
RN 651330-64-2 HCAPLUS

CN 2-Piperidinecarboxylic acid, 3-[[[4-[[[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]amino]methyl]phenyl]methyl]amino]propyl ester (9CI) (CA INDEX NAME)

PAGE 1-A

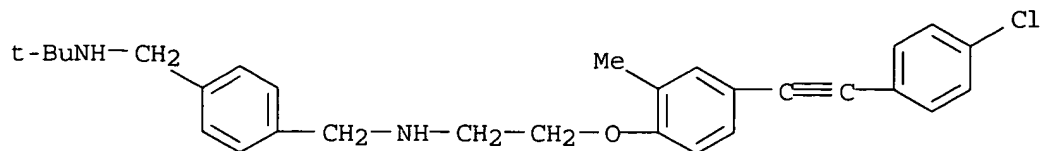


PAGE 1-B



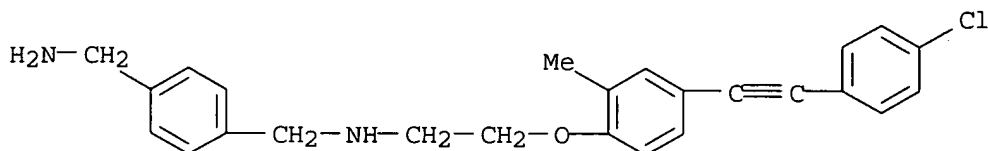
RN 651330-65-3 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-(1,1-dimethylethyl)- (9CI) (CA INDEX NAME)



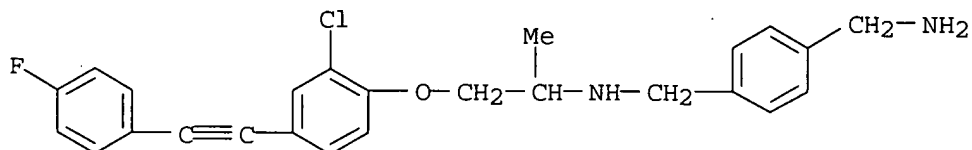
RN 651330-68-6 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]- (9CI) (CA INDEX NAME)



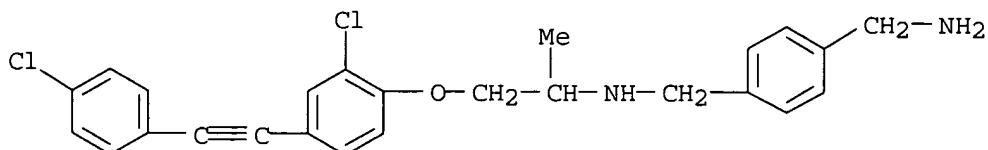
RN 773847-69-1 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[2-chloro-4-[(4-fluorophenyl)ethynyl]phenoxy]-1-methylethyl]- (9CI) (CA INDEX NAME)



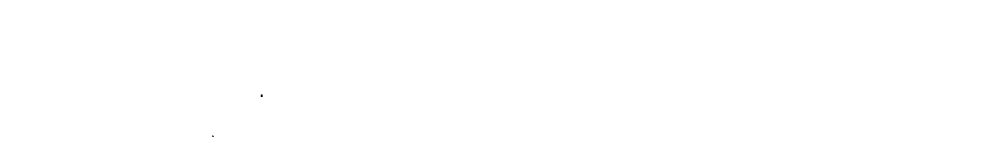
RN 776293-45-9 HCAPLUS

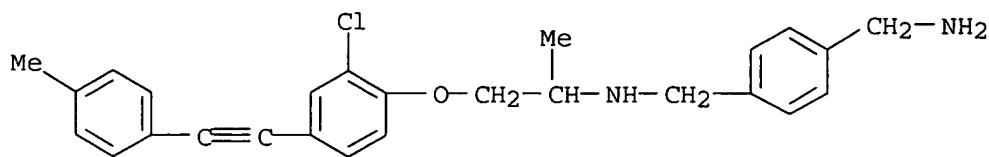
CN 1,4-Benzenedimethanamine, N-[2-[2-chloro-4-[(4-chlorophenyl)ethynyl]phenoxy]-1-methylethyl]- (9CI) (CA INDEX NAME)



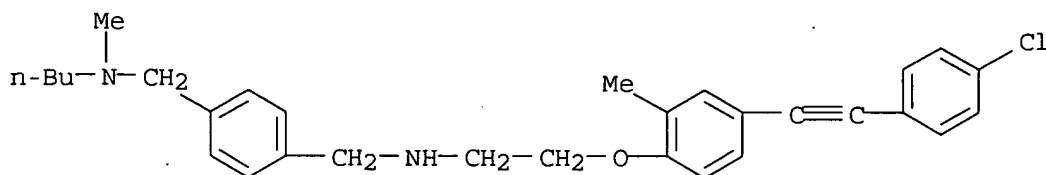
RN 792905-81-8 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[2-chloro-4-[(4-methylphenyl)ethynyl]phenoxy]-1-methylethyl]- (9CI) (CA INDEX NAME)

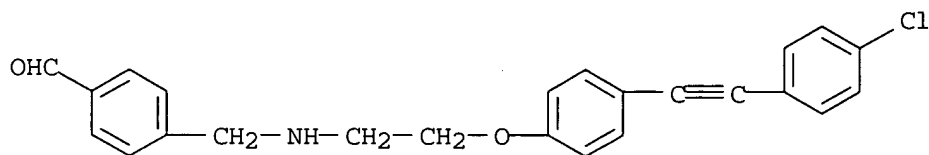




RN 832116-38-8 HCAPLUS  
 CN 1,4-Benzenedimethanamine, N-butyl-N'-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N-methyl- (9CI) (CA INDEX NAME)



IT 651330-74-4P  
 RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
 (preparation of acetylenic compound useful in the treatment of inflammatory disorders)  
 RN 651330-74-4 HCAPLUS  
 CN Benzaldehyde, 4-[[[2-[4-[(4-chlorophenyl)ethynyl]phenoxy]ethyl]amino]methyl]- (9CI) (CA INDEX NAME)



L4 ANSWER 2 OF 4 HCAPLUS COPYRIGHT 2005 ACS on STN  
 ACCESSION NUMBER: 2004:650899 HCAPLUS  
 DOCUMENT NUMBER: 141:173978  
 TITLE: Preparation of aminoacetonitrile derivatives as agricultural and horticultural insecticides  
 INVENTOR(S): Andoh, Nobuharu; Sanpei, Osamu; Sakata, Kazuyuki  
 PATENT ASSIGNEE(S): Nihon Nohyaku Co., Ltd., Japan  
 SOURCE: Eur. Pat. Appl., 48 pp.  
 CODEN: EPXXDW  
 DOCUMENT TYPE: Patent  
 LANGUAGE: English  
 FAMILY ACC. NUM. COUNT: 2  
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 1445251	A1	20040811	EP 2004-10346	19990428
R: CH, DE, FR, GB, IT, LI				
EP 953565	A2	1999-1-103	EP 1999-107461	19990428
EP 953565	A3	20021204		



EP 953565 B1 20040908

R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,  
IE, SI, LT, LV, FI, RO

PRIORITY APPLN. INFO.:

JP 1998-137806

A 19980501

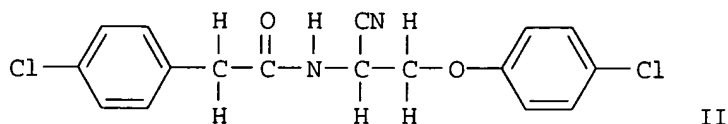
EP 1999-107461

A3 19990428

OTHER SOURCE(S):

MARPAT 141:173978

GI



AB The title compds. Ar1(Q)dC(O)NR3C(CN)R4(CR5R6)aW(CR7R8)bAr2 [I; Ar1, Ar2 = (substituted) Ph, (substituted) phenyloxy, (substituted) phenylacetylene; (substituted) pyridyl and (substituted) naphthyl; Q = CR1R2 (wherein R1, R2 = H, halo, (halo)alkyl, etc.); R3 = H, (halo)alkyl, etc.; R4-R8 = H, halo, (halo)alkyl, etc.; W = O, S, SO2 or NR9 (wherein R9 = H, alkyl); a, b = 0-4; d = 0-1], useful as insecticides, were prepared E.g., a multi-step synthesis of II (starting from 4-chlorophenol and bromoacetaldehyde dimethylacetal), was given. The compds. I were tested against diamondback moth and against smaller tea tortrix (data were given for representative compds. I).

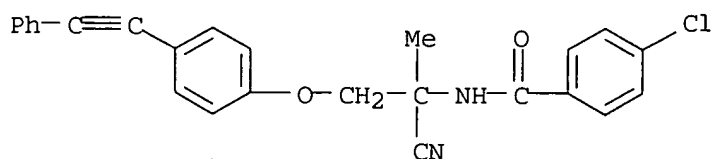
IT 736172-93-3P

RL: AGR (Agricultural use); BSU (Biological study, unclassified); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of aminoacetonitrile derivs. as agricultural and horticultural insecticides)

RN 736172-93-3 HCAPLUS

CN Benzamide, 4-chloro-N-[1-cyano-1-methyl-2-[4-(phenylethynyl)phenoxy]ethyl]-(9CI) (CA INDEX NAME)



L4 ANSWER 3 OF 4 HCAPLUS COPYRIGHT 2005 ACS on STN

ACCESSION NUMBER: 2004:80636 HCAPLUS

DOCUMENT NUMBER: 140:145889

TITLE: Preparation of acetylenic-aryl-methylamines useful in treating inflammatory disorders

INVENTOR(S): Beers, Scott; Mailloy, Elizabeth A.; Wachter, Michael P.

PATENT ASSIGNEE(S): Janssen Pharmaceutica N.V., Belg.

SOURCE: PCT Int. Appl., 61 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

## PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004009531	A1	20040129	WO 2003-US23140	20030724
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
PRIORITY APPLN. INFO.:			US 2002-398138P	P 20020724
OTHER SOURCE(S):		MARPAT 140:145889		
GI				

\* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT \*

AB Title compds. I [A = cycloalkyldiyl, cyclic heteroalkyldiyl, etc.; B, E = (hetero)aryldiyl; R1 = cycloalkyl, cyclic heteroalkyl, heteroaryl, etc.; R2 = H, alkanyl, alkoxy, acyl, etc.; R3-4 = H, alkanyl, carboxy, cycloalkyl, etc.] are prepared For instance, 4-iodo-2-methylphenol is alkylated with chloroacetone (acetone, K<sub>2</sub>CO<sub>3</sub>, NaI, reflux, 18 h), the product coupled to 4-chlorophenylacetylene (Pd(II), Et<sub>3</sub>N) and used to alkylate p-xylylene diamine (CH<sub>2</sub>Cl<sub>2</sub>, HOAc, NaBH(OAc)<sub>3</sub>) to give II. In an NADPH oxidase assay for inhibition of superoxide-mediated reduction of cytochrome C in human neutrophils incubated with phorbol myristate acetate, 26 compds. I had IC<sub>50</sub> values of 0.7-11.13 µM; II had IC<sub>50</sub> = 0.7 µM. I are useful in treating or ameliorating reactive oxygen species-mediated inflammatory disorders.

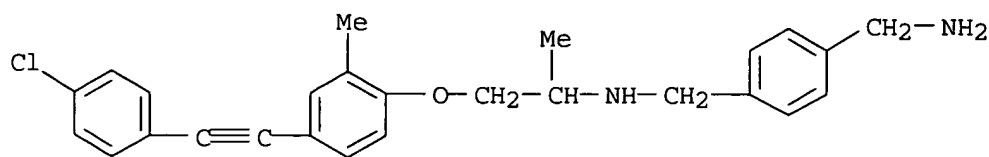
IT 651330-36-8P 651330-37-9P 651330-38-0P  
651330-39-1P 651330-40-4P 651330-41-5P  
651330-42-6P 651330-43-7P 651330-44-8P  
651330-45-9P 651330-46-0P 651330-47-1P  
651330-48-2P 651330-49-3P 651330-50-6P  
651330-51-7P 651330-52-8P 651330-53-9P  
651330-54-0P 651330-55-1P 651330-57-3P  
651330-58-4P 651330-59-5P 651330-61-9P  
651330-63-1P 651330-64-2P 651330-65-3P  
651330-66-4P 651330-68-6P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

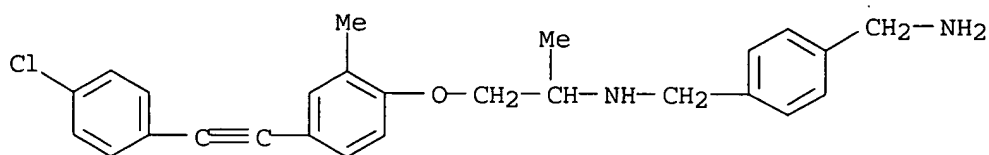
(preparation of acetylenic-aryl-methylamines as NAD oxidase hydride donor inhibitors useful in treating reactive oxygen species-mediated inflammatory disorders)

RN 651330-36-8 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]-1-methylethyl]- (9CI) (CA INDEX NAME)

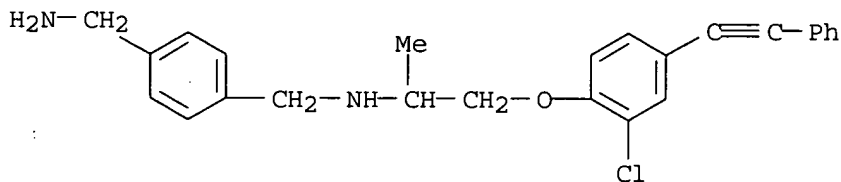


RN 651330-37-9 HCAPLUS  
 CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]-1-methylethyl]-, dihydrochloride (9CI) (CA INDEX NAME)

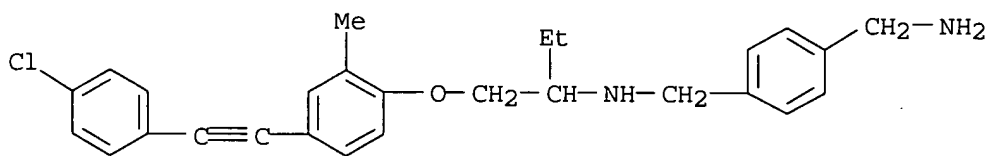


● 2 HCl

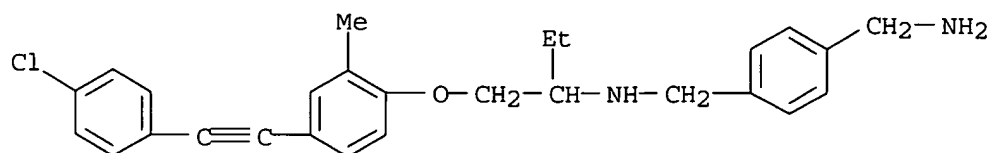
RN 651330-38-0 HCAPLUS  
 CN 1,4-Benzenedimethanamine, N-[2-[2-chloro-4-(phenylethynyl)phenoxy]-1-methylethyl]- (9CI) (CA INDEX NAME)



RN 651330-39-1 HCAPLUS  
 CN 1,4-Benzenedimethanamine, N-[1-[[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]methyl]propyl]- (9CI) (CA INDEX NAME)



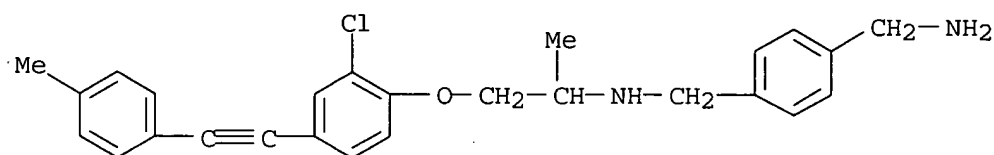
RN 651330-40-4 HCAPLUS  
 CN 1,4-Benzenedimethanamine, N-[1-[[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]methyl]propyl]-, dihydrochloride (9CI) (CA INDEX NAME)



● 2 HCl

RN 651330-41-5 HCAPLUS

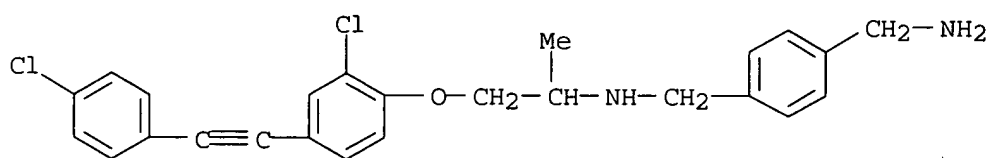
CN 1,4-Benzenedimethanamine, N-[2-[2-chloro-4-[(4-methylphenyl)ethynyl]phenoxy]-1-methylethyl]-, dihydrochloride (9CI) (CA INDEX NAME)



● 2 HCl

RN 651330-42-6 HCAPLUS

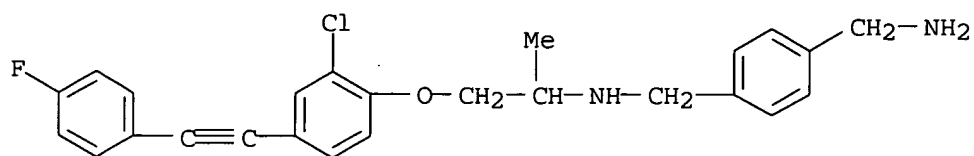
CN 1,4-Benzenedimethanamine, N-[2-[2-chloro-4-[(4-chlorophenyl)ethynyl]phenoxy]-1-methylethyl]-, dihydrochloride (9CI) (CA INDEX NAME)



● 2 HCl

RN 651330-43-7 HCAPLUS

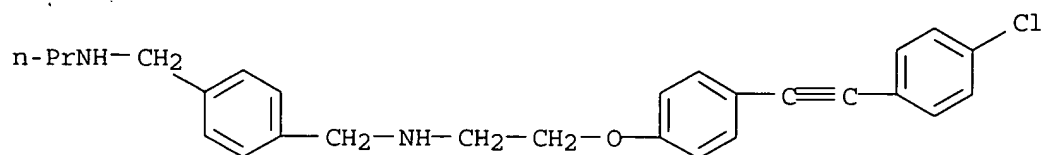
CN 1,4-Benzenedimethanamine, N-[2-[2-chloro-4-[(4-fluorophenyl)ethynyl]phenoxy]-1-methylethyl]-, dihydrochloride (9CI) (CA INDEX NAME)



●2 HCl

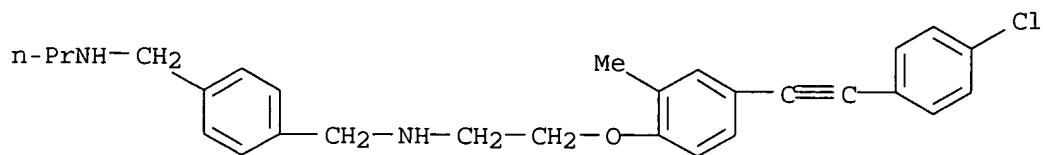
RN 651330-44-8 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]phenoxy]ethyl]-N'-propyl- (9CI) (CA INDEX NAME)



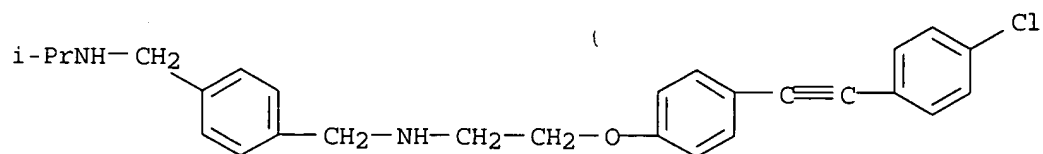
RN 651330-45-9 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-propyl- (9CI) (CA INDEX NAME)



RN 651330-46-0 HCAPLUS

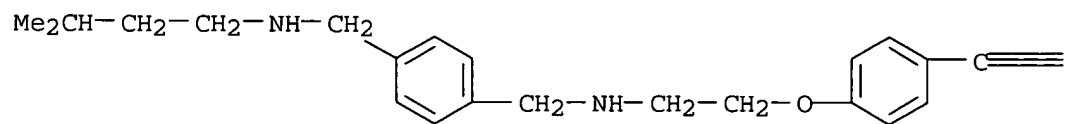
CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]phenoxy]ethyl]-N'-(1-methylethyl)- (9CI) (CA INDEX NAME)



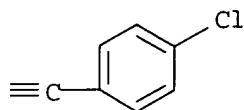
RN 651330-47-1 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]phenoxy]ethyl]-N'-(3-methylbutyl)- (9CI) (CA INDEX NAME)

PAGE 1-A



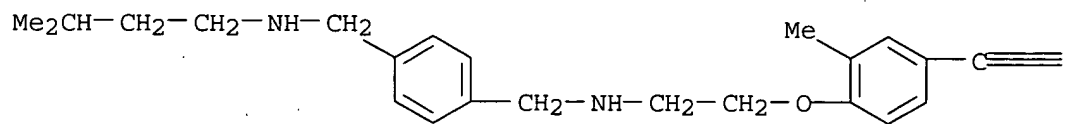
PAGE 1-B



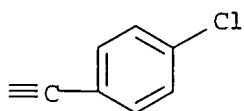
RN 651330-48-2 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-(3-methylbutyl)- (9CI) (CA INDEX NAME)

PAGE 1-A



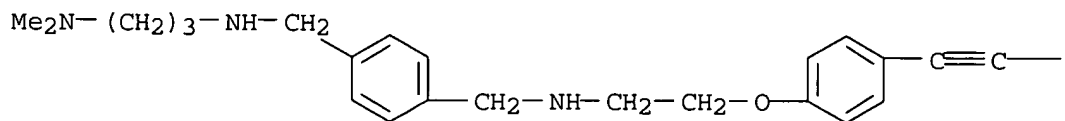
PAGE 1-B



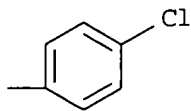
RN 651330-49-3 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]phenoxy]ethyl]-N'-[3-(dimethylamino)propyl]- (9CI) (CA INDEX NAME)

PAGE 1-A



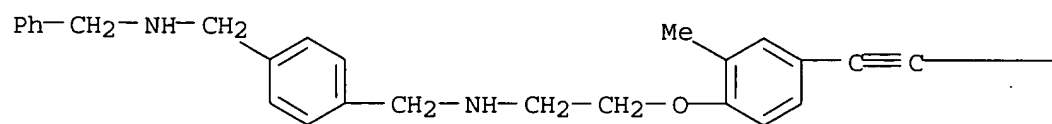
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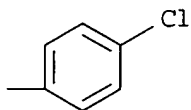
RN 651330-50-6 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-(phenylmethyl)- (9CI) (CA INDEX NAME)

PAGE 1-A



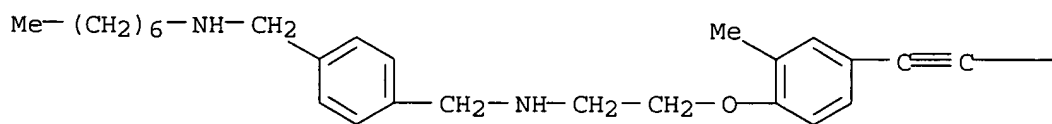
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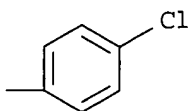
RN 651330-51-7 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-heptyl- (9CI) (CA INDEX NAME)

PAGE 1-A



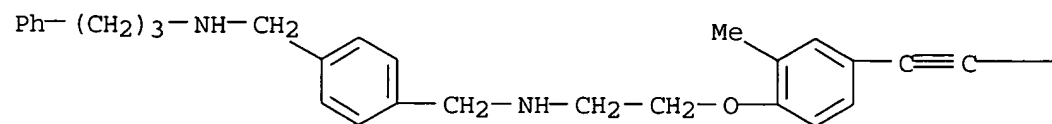
PAGE 1-B



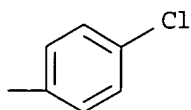
RN 651330-52-8 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-(3-phenylpropyl)- (9CI) (CA INDEX NAME)

PAGE 1-A



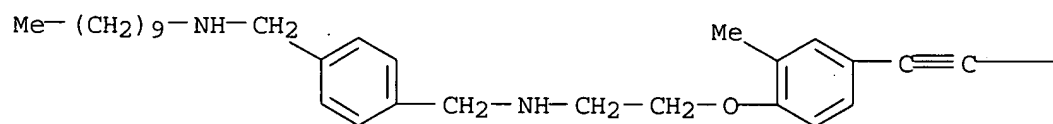
PAGE 1-B



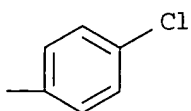
RN 651330-53-9 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-decyl- (9CI) (CA INDEX NAME)

PAGE 1-A



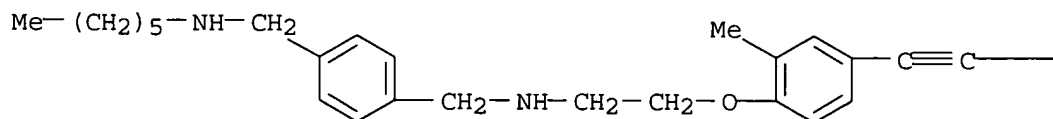
PAGE 1-B



RN 651330-54-0 HCAPLUS

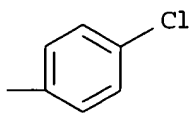
CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-hexyl- (9CI) (CA INDEX NAME)

PAGE 1-A





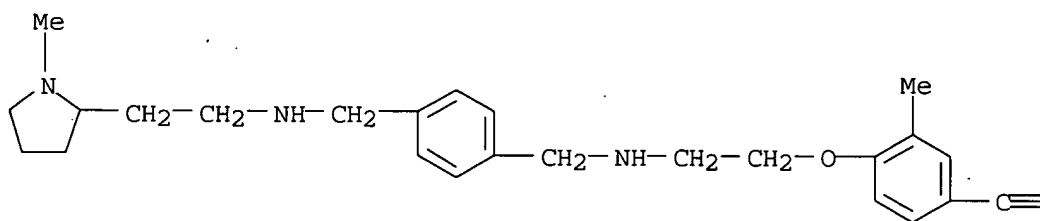
PAGE 1-B



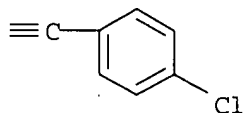
RN 651330-55-1 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-[2-(1-methyl-2-pyrrolidinyl)ethyl]- (9CI) (CA INDEX NAME)

PAGE 1-A



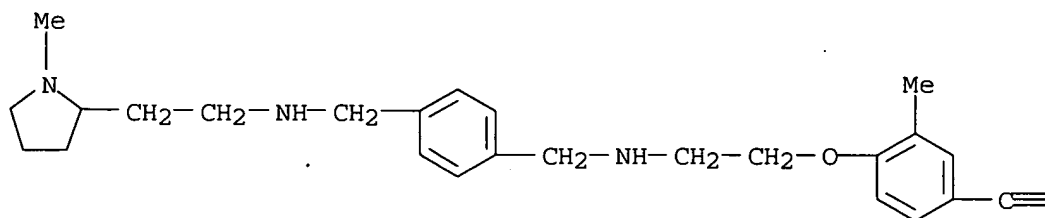
PAGE 1-B



RN 651330-57-3 HCAPLUS

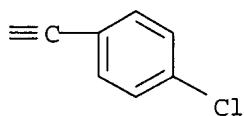
CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-[2-(1-methyl-2-pyrrolidinyl)ethyl]-, trihydrochloride (9CI) (CA INDEX NAME)

PAGE 1-A



● 3 HCl

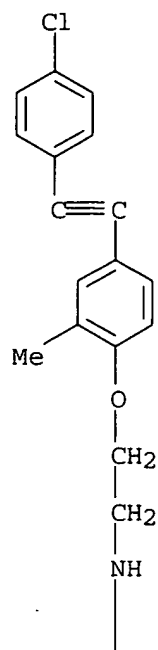
PAGE 1-B



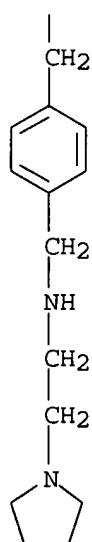
RN 651330-58-4 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-[2-(1-pyrrolidinyl)ethyl]- (9CI) (CA INDEX NAME)

PAGE 1-A



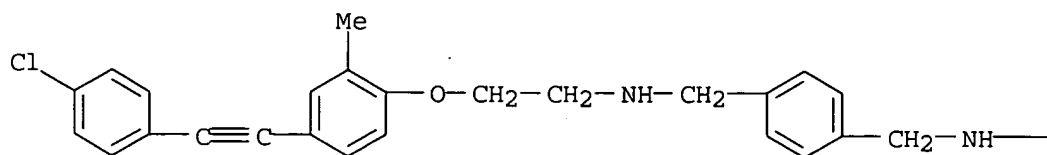
PAGE 2-A



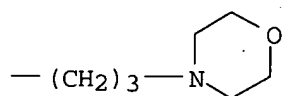
RN 651330-59-5 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-[3-(4-morpholinyl)propyl]- (9CI) (CA INDEX NAME)

PAGE 1-A

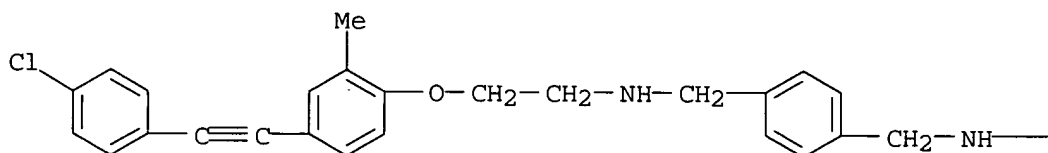


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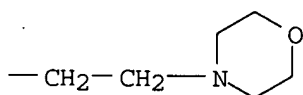


RN 651330-61-9 HCAPLUS  
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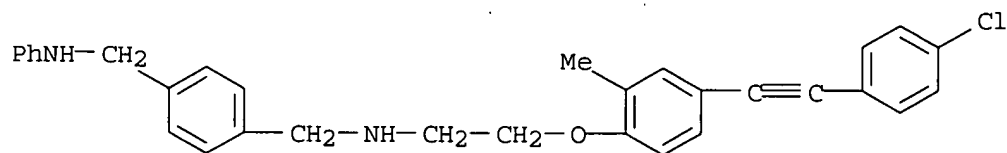
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PAGE 1-B



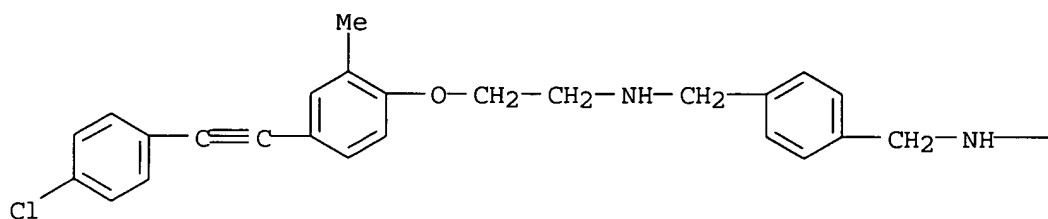
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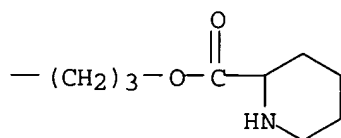
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 CN 2-Piperidinecarboxylic acid, 3-[[[4-[[[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]amino]methyl]phenyl]methyl]amino]propyl ester (9CI)

(CA INDEX NAME)

PAGE 1-A

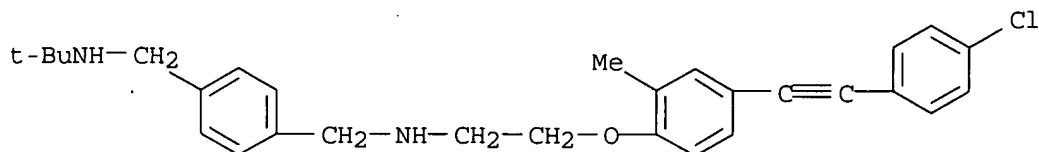


PAGE 1-B



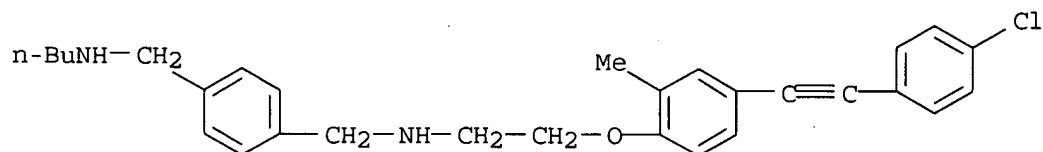
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CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]-N'-(1,1-dimethylethyl)- (9CI) (CA INDEX NAME)



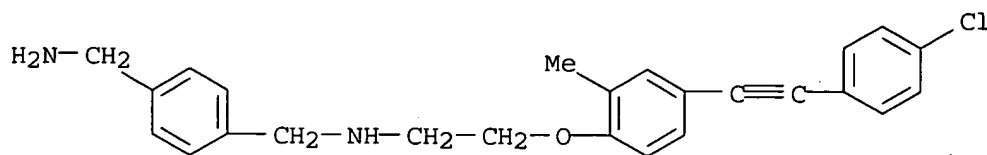
RN 651330-66-4 HCAPLUS

CN 1,4-Benzenedimethanamine, N-butyl-N'-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]- (9CI) (CA INDEX NAME)



RN 651330-68-6 HCAPLUS

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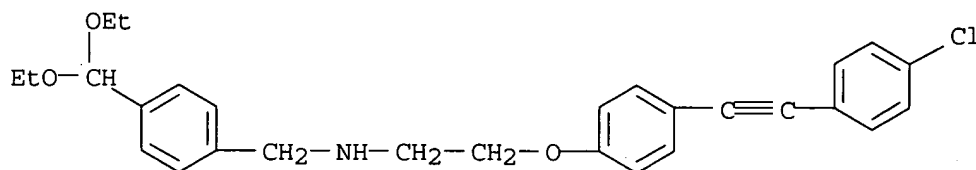
IT 651330-73-3P 651330-74-4P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of acetylenic-aryl-methylamines as NAD oxidase hydride donor inhibitors useful in treating reactive oxygen species-mediated inflammatory disorders)

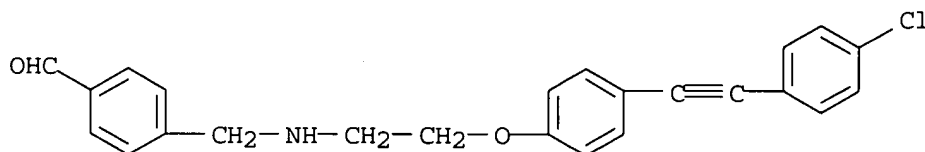
RN 651330-73-3 HCAPLUS

CN Benzenemethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]phenoxy]ethyl]-4-(diethoxymethyl)- (9CI) (CA INDEX NAME)



RN 651330-74-4 HCAPLUS

CN Benzaldehyde, 4-[[[2-[4-[(4-chlorophenyl)ethynyl]phenoxy]ethyl]amino]methyl]- (9CI) (CA INDEX NAME)



REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 4 OF 4 HCAPLUS COPYRIGHT 2005 ACS on STN

ACCESSION NUMBER: 1999:708444 HCAPLUS

DOCUMENT NUMBER: 131:310455

TITLE: Preparation of aroylaminoacetonitriles as agricultural and horticultural insecticides

INVENTOR(S): Andoh, Nobuharu; Sanpei, Osamu; Sakata, Kazuyuki

PATENT ASSIGNEE(S): Nihon Nohyaku Co., Ltd., Japan

SOURCE: Eur. Pat. Appl., 63 pp.

CODEN: EPXXDW

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 2

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 953565	A2	19991103	EP 1999-107461	19990428

EP 953565 A3 20021204  
 EP 953565 B1 20040908  
 R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,  
 IE, SI, LT, LV, FI, RO  
US 6239077 B1 20010529 US 1999-295319 19990421  
 TW 585849 B 20040501 TW 1999-88106732 19990427  
 EP 1445251 A1 20040811 EP 2004-10346 19990428  
 R: CH, DE, FR, GB, IT, LI  
 CN 1234177 A 19991110 CN 1999-105289 19990430  
 CN 1132516 B 20031231  
 AU 9926027 A1 19991111 AU 1999-26027 19990430  
 AU 752112 B2 20020905  
 JP 2000026392 A2 20000125 JP 1999-124560 19990430  
 PRIORITY APPLN. INFO.: JP 1998-137806 A 19980501  
 EP 1999-107461 A3 19990428

OTHER SOURCE(S): MARPAT 131:310455

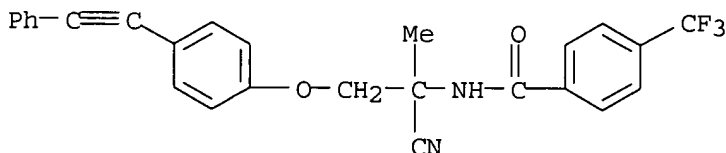
AB Ar1QdCONR3C(CN)R4(CR5R6)aW(CR7R8)bAr2 [I; Ar1, Ar2 = (substituted) Ph, PhO, pyridyl, pyridyloxy, naphthyl; Q = CR1R2; R1, R2 = H, halo, (halo)alkyl, (halo)alkoxy, (substituted) cycloalkyl; R1R2 = (substituted) C2-6 alkylene, CH:CH, C.tplbond.C; d = 0, 1; R3 = H, (halo)alkyl; R4-R8 = H, halo, (halo)alkyl; W = O, S, SO2, NR9; R9 = H, alkyl; a, b = 0-4], were prepared. Thus, 4-chlorophenol, bromoacetaldehyde di-Me acetal, K2CO3, and cat. NaI were refluxed 3 h in DMF to give 4-chlorophenoxyacetaldehyde di-Me acetal. This was refluxed with aqueous HCl in acetone to give crude 4-chlorophenoxyacetaldehyde, which was stirred with NaCN and NH4Cl in aqueous NH3 to give a residue. This was stirred with 4-chlorophenylacetyl chloride and Et3N in THF to give I (Ar1, Ar2 = 4-ClC6H4; R1-R8 = H; W = O; a, d = 1; b = 0). Numerous I at 500 ppm gave 100% kill of Plutella xylostella on cabbage seedlings.

IT 247199-30-0P

RL: AGR (Agricultural use); BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); USES (Uses) (preparation of aroylaminoacetonitriles as agricultural and horticultural insecticides)

RN 247199-30-0 HCAPLUS

CN Benzamide, N-[1-cyano-1-methyl-2-[4-(phenylethynyl)phenoxy]ethyl]-4-(trifluoromethyl)- (9CI) (CA INDEX NAME)



=> FIL REGISTRY

COST IN U.S. DOLLARS

SINCE FILE ENTRY	TOTAL SESSION
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FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE ENTRY	TOTAL SESSION
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CA SUBSCRIBER PRICE

FILE 'REGISTRY' ENTERED AT 08:43:44 ON 30 AUG 2005

08/30/2005 10626155.trn

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STRUCTURE FILE UPDATES: 29 AUG 2005 HIGHEST RN 862072-85-3  
DICTIONARY FILE UPDATES: 29 AUG 2005 HIGHEST RN 862072-85-3

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TSCA INFORMATION NOW CURRENT THROUGH JULY 14, 2005

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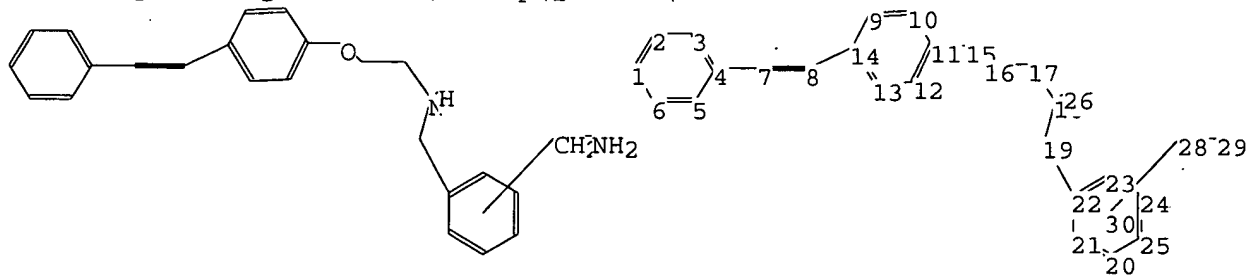
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\* The CA roles and document type information have been removed from \*  
\* the IDE default display format and the ED field has been added, \*  
\* effective March 20, 2005. A new display format, IDERL, is now \*  
\* available and contains the CA role and document type information. \*  
\*  
\*\*\*\*\*

Structure search iteration limits have been increased. See HELP SLIMITS  
for details.

Experimental and calculated property data are now available. For more  
information enter HELP PROP at an arrow prompt in the file or refer  
to the file summary sheet on the web at:  
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>

Uploading C:\Program Files\Stnexp\Queries\10626155a.str



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ring nodes :

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chain bonds :



08/30/2005 10626155.trn

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11-15 15-16 17-18 18-19  
exact bonds :  
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containing 1 : 9 : 20 :

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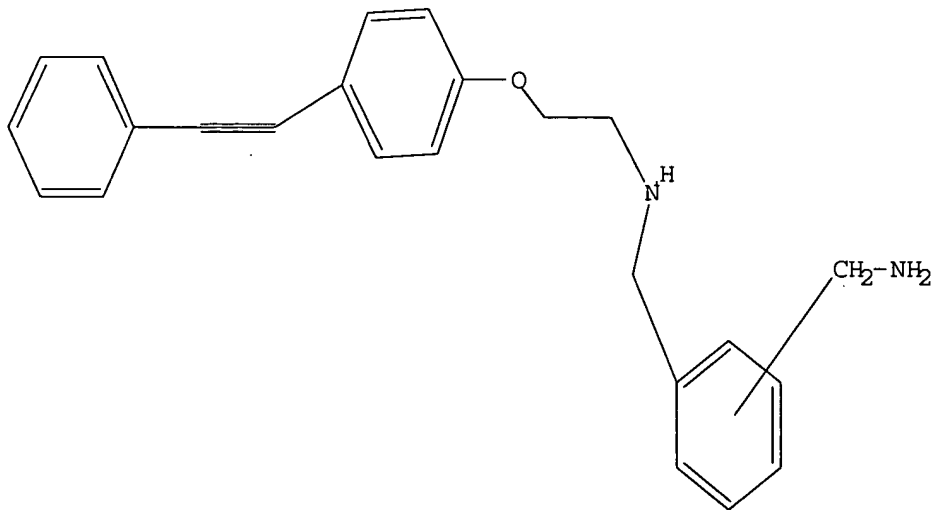
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19:CLASS 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom 26:CLASS 28:CLASS  
29:CLASS 30:CLASS

L5 STRUCTURE UPLOADED

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L5 HAS NO ANSWERS

L5 STR



Structure attributes must be viewed using STN Express query preparation.

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SAMPLE SCREEN SEARCH COMPLETED - 9 TO ITERATE

100.0% PROCESSED

9 ITERATIONS

0 ANSWERS

08/30/2005 10626155.trn

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
PROJECTED ITERATIONS: 9 TO 360  
PROJECTED ANSWERS: 0 TO 0

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=> s l5 sss full  
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FULL SCREEN SEARCH COMPLETED - 327 TO ITERATE

100.0% PROCESSED 327 ITERATIONS  
SEARCH TIME: 00.00.01

12 ANSWERS

L7 12 SEA SSS FUL L5

=> FIL HCAPLUS  
COST IN U.S. DOLLARS

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ENTRY	SESSION
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FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
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FILE 'HCAPLUS' ENTERED AT 08:44:18 ON 30 AUG 2005  
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FILE COVERS 1907 - 30 Aug 2005 VOL 143 ISS 10  
FILE LAST UPDATED: 29 Aug 2005 (20050829/ED)

New CAS Information Use Policies, enter HELP USAGETERMS for details.

This file contains CAS Registry Numbers for easy and accurate substance identification.

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L8

2 L7

=> d l8 ibib abs hitstr tot

L8 ANSWER 1 OF 2 HCAPLUS COPYRIGHT 2005 ACS on STN  
ACCESSION NUMBER: 2005:78301 HCAPLUS  
DOCUMENT NUMBER: 142:176433

08/30/2005 10626155.trn

TITLE: A preparation of acetylenic compounds, useful in the treatment of inflammatory disorders

INVENTOR(S): Beers, Scott; Malloy, Elizabeth A.; Wachter, Michael P.

PATENT ASSIGNEE(S): USA

SOURCE: U.S. Pat. Appl. Publ., 19 pp.  
CODEN: USXXCO

DOCUMENT TYPE: Patent

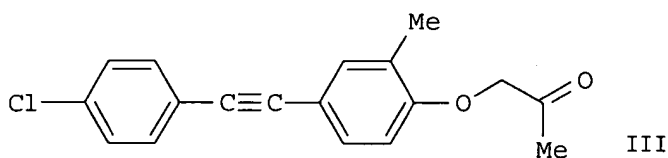
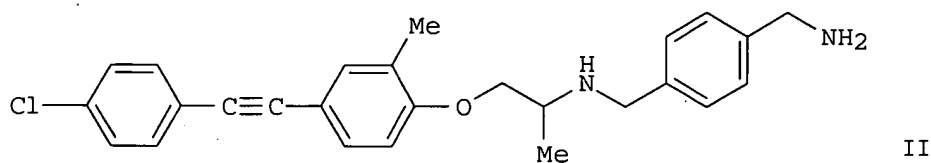
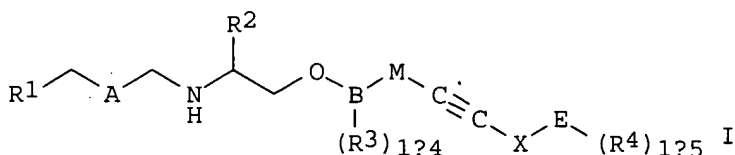
LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 2005020838	A1	20050127	US 2003-626155	20030724
PRIORITY APPLN. INFO.:			US 2003-626155	20030724
OTHER SOURCE(S):	MARPAT 142:176433			

GI



AB The invention relates to a preparation of acetylenic compds. of formula I [wherein: A is cycloalkyldiyl, cyclic heteroalkyldiyl, or (hetero)aryldiyl; B is (hetero)aryldiyl; E is (hetero)aryldiyl; M and X are independently (CH<sub>2</sub>)<sub>0-4</sub>; R<sub>1</sub> is cycloalkyl, cyclic heteroalkyl, or (hetero)aryl, etc.; R<sub>2</sub> is H, alkyl, alkoxy, CHO, CO<sub>2</sub>H, or NH<sub>2</sub>, etc.; R<sub>3</sub> and R<sub>4</sub> are independently H, alkyl, CHO, cycloalkyl, or aryl, etc.], useful in the treatment of inflammatory disorders. For instance, acetylene derivative II (oxidase inhibition: IC<sub>50</sub> = 0.7 μM) was prepared via reductive amination of ketone III by 1,4-bis(aminomethyl)benzene.

IT 651330-36-8P 651330-37-9P 651330-38-0P  
651330-39-1P 651330-68-6P 773847-69-1P  
776293-45-9P 792905-81-8P

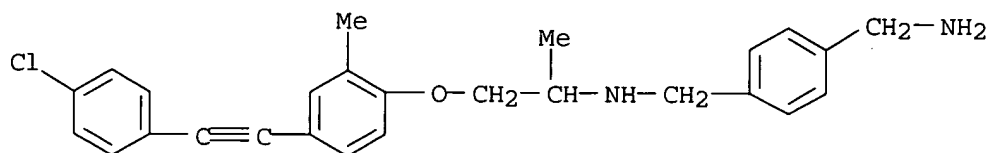
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU

(Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of acetylenic compound useful in the treatment of inflammatory disorders)

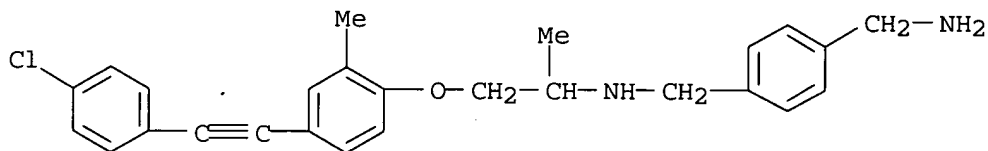
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CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]-1-methylethyl]- (9CI) (CA INDEX NAME)



RN 651330-37-9 HCAPLUS

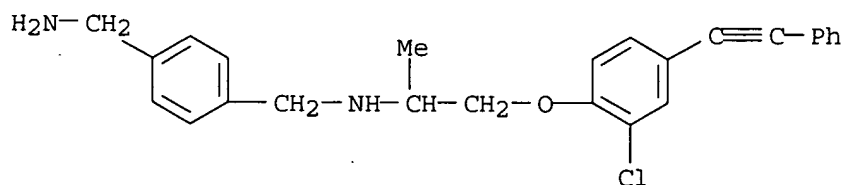
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● 2 HCl

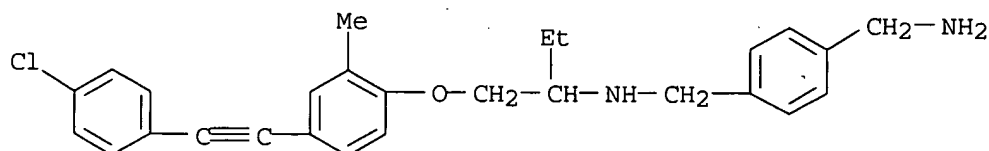
RN 651330-38-0 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[2-chloro-4-(phenylethynyl)phenoxy]-1-methylethyl]- (9CI) (CA INDEX NAME)



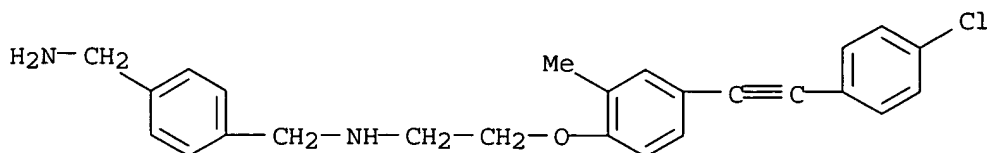
RN 651330-39-1 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[1-[[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]methyl]propyl]- (9CI) (CA INDEX NAME)



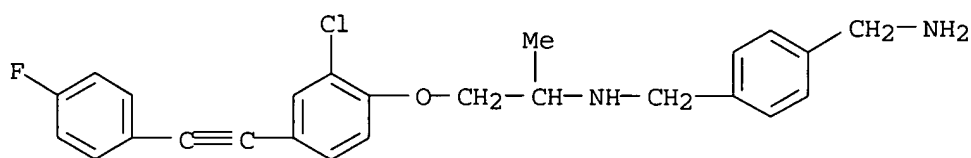
RN 651330-68-6 HCAPLUS

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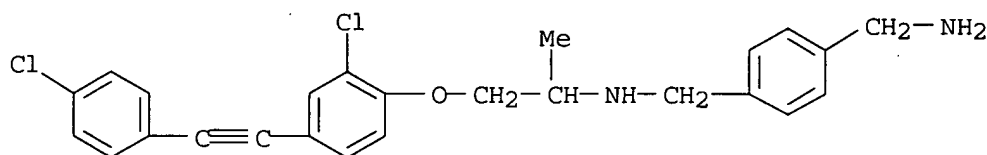
RN 773847-69-1 HCAPLUS

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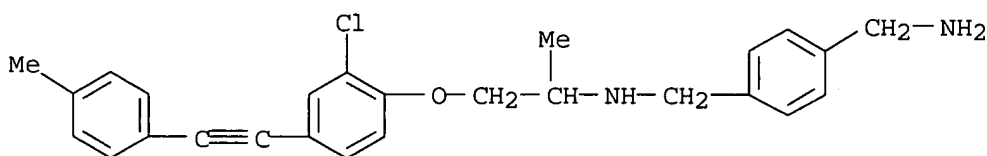
RN 776293-45-9 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[2-chloro-4-[(4-chlorophenyl)ethynyl]phenoxy]-1-methylethyl] - (9CI) (CA INDEX NAME)



RN 792905-81-8 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[2-chloro-4-[(4-methylphenyl)ethynyl]phenoxy]-1-methylethyl] - (9CI) (CA INDEX NAME)



L8 ANSWER 2 OF 2 HCAPLUS COPYRIGHT 2005 ACS on STN

ACCESSION NUMBER: 2004:80636 HCAPLUS

DOCUMENT NUMBER: 140:145889

TITLE: Preparation of acetylenic-aryl-methylamines useful in treating inflammatory disorders

INVENTOR(S): Beers, Scott; Malloy, Elizabeth A.; Wachter, Michael P.

PATENT ASSIGNEE(S): Janssen Pharmaceutica N.V., Belg.

SOURCE: PCT Int. Appl., 61 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English  
 FAMILY ACC. NUM. COUNT: 1  
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004009531	A1	20040129	WO 2003-US23140	20030724
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
PRIORITY APPLN. INFO.:			US 2002-398138P	P. 20020724
OTHER SOURCE(S):			MARPAT 140:145889	
GI				

\* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT \*

AB Title compds. I [A = cycloalkyldiyl, cyclic heteroalkyldiyl, etc.; B, E = (hetero)aryldiyl; R1 = cycloalkyl, cyclic heteroalkyl, heteroaryl, etc.; R2 = H, alkanyl, alkoxy, acyl, etc.; R3-4 = H, alkanyl, carboxy, cycloalkyl, etc.] are prepared For instance, 4-iodo-2-methylphenol is alkylated with chloroacetone (acetone, K<sub>2</sub>CO<sub>3</sub>, NaI, reflux, 18 h), the product coupled to 4-chlorophenylacetylene (Pd(II), Et<sub>3</sub>N) and used to alkylate p-xylylene diamine (CH<sub>2</sub>Cl<sub>2</sub>, HOAc, NaBH(OAc)<sub>3</sub>) to give II. In an NADPH oxidase assay for inhibition of superoxide-mediated reduction of cytochrome C in human neutrophils incubated with phorbol myristate acetate, 26 compds. I had IC<sub>50</sub> values of 0.7-11.13 µM; II had IC<sub>50</sub> = 0.7 µM. I are useful in treating or ameliorating reactive oxygen species-mediated inflammatory disorders.

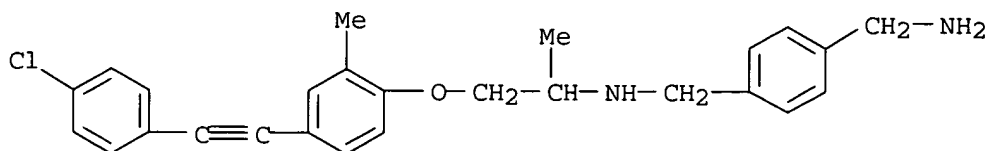
IT 651330-36-8P 651330-37-9P 651330-38-0P  
 651330-39-1P 651330-40-4P 651330-41-5P  
 651330-42-6P 651330-43-7P 651330-68-6P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of acetylenic-aryl-methylamines as NAD oxidase hydride donor inhibitors useful in treating reactive oxygen species-mediated inflammatory disorders)

RN 651330-36-8 HCAPLUS

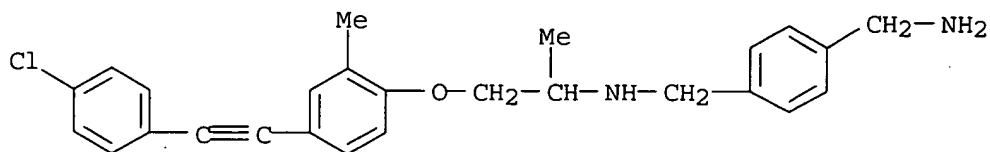
CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]-1-methylethyl]- (9CI) (CA INDEX NAME)



08/30/2005 10626155.trn

RN 651330-37-9 HCAPLUS

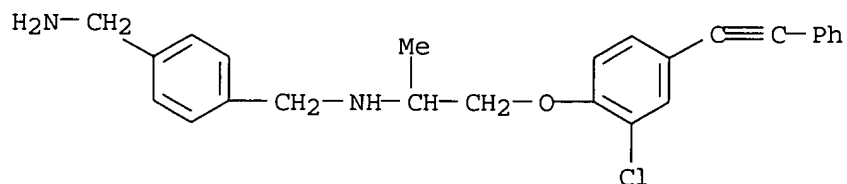
CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]-1-methylethyl]-, dihydrochloride (9CI) (CA INDEX NAME)



● 2 HCl

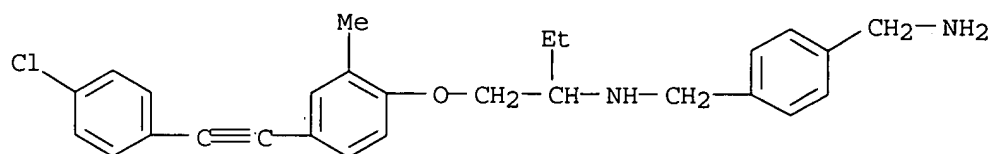
RN 651330-38-0 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[2-chloro-4-(phenylethynyl)phenoxy]-1-methylethyl]- (9CI) (CA INDEX NAME)



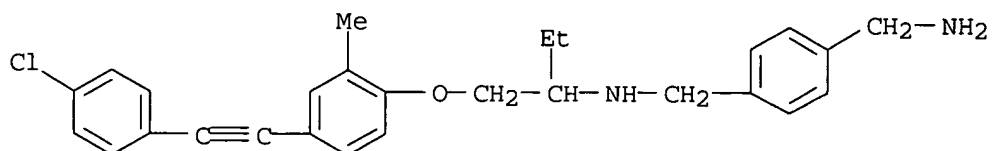
RN 651330-39-1 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[1-[[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]methyl]propyl]- (9CI) (CA INDEX NAME)



RN 651330-40-4 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[1-[[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]methyl]propyl]-, dihydrochloride (9CI) (CA INDEX NAME)

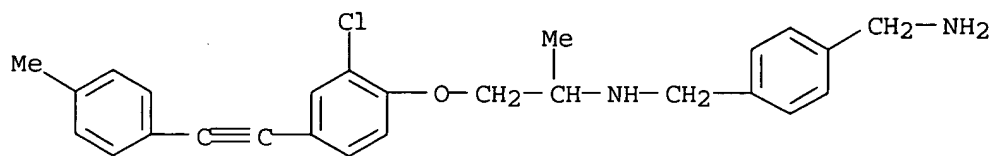


● 2 HCl

08/30/2005 10626155.trn

RN 651330-41-5 HCAPLUS

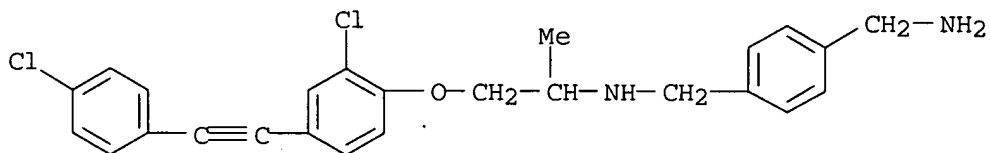
CN 1,4-Benzenedimethanamine, N-[2-[2-chloro-4-[(4-methylphenyl)ethynyl]phenoxy]-1-methylethyl]-, dihydrochloride (9CI) (CA INDEX NAME)



● 2 HCl

RN 651330-42-6 HCAPLUS

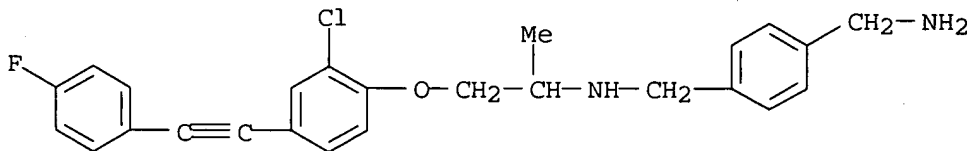
CN 1,4-Benzenedimethanamine, N-[2-[2-chloro-4-[(4-chlorophenyl)ethynyl]phenoxy]-1-methylethyl]-, dihydrochloride (9CI) (CA INDEX NAME)



● 2 HCl

RN 651330-43-7 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[2-chloro-4-[(4-fluorophenyl)ethynyl]phenoxy]-1-methylethyl]-, dihydrochloride (9CI) (CA INDEX NAME)

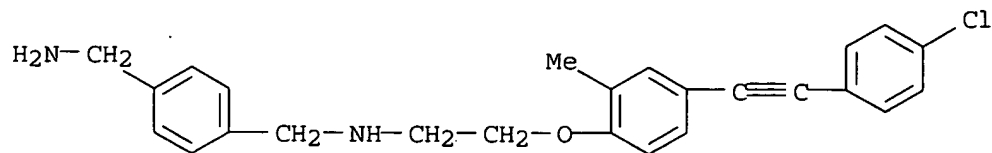


● 2 HCl

RN 651330-68-6 HCAPLUS

CN 1,4-Benzenedimethanamine, N-[2-[4-[(4-chlorophenyl)ethynyl]-2-methylphenoxy]ethyl]- (9CI) (CA INDEX NAME)





REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
12.33	379.46

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
-1.46	-4.38

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